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Size and symmetry of the superconducting gap in the f.c.c.  $\text{Cs}_3\text{C}_{60}$  polymorph close to the metal-Mott insulator boundary

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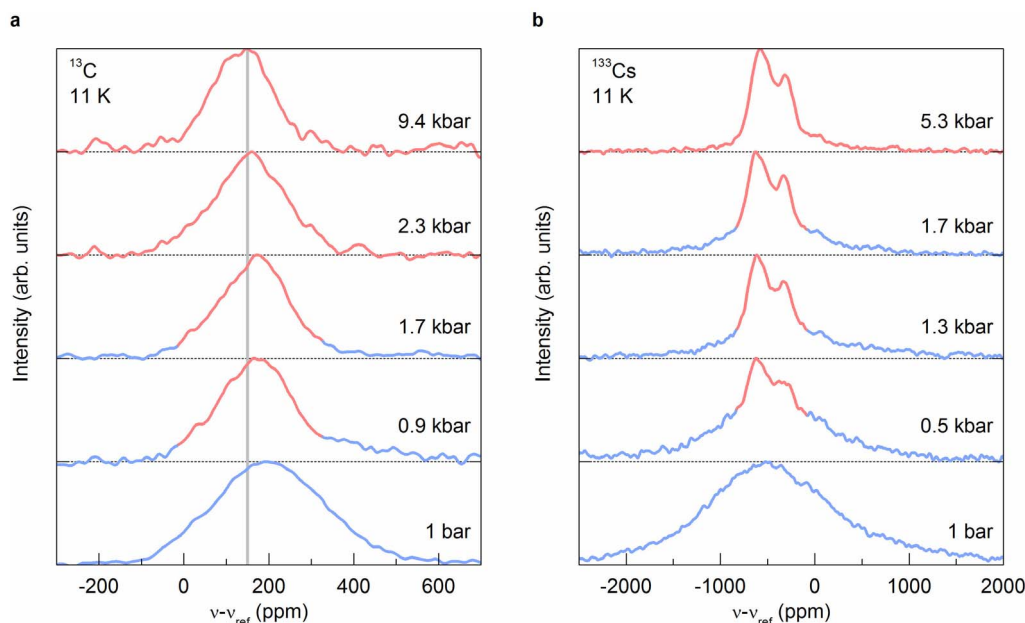
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The alkali fullerenes,  $\text{A}_3\text{C}_{60}$  ( $\text{A}$  = alkali metal) are molecular superconductors that undergo a transition to a magnetic Mott-insulating state at large lattice parameters. However, although the size and the symmetry of the superconducting gap,  $\Delta$ , are both crucial for the understanding of the pairing mechanism, they are currently unknown for superconducting fullerenes close to the correlation-driven magnetic insulator. Here we report a comprehensive nuclear magnetic resonance (NMR) study of face-centred-cubic (f.c.c.)  $\text{Cs}_3\text{C}_{60}$  polymorph, which can be tuned continuously through the bandwidth-controlled Mott insulator-metal/superconductor transition by pressure. When superconductivity emerges from the insulating state at large interfullerene separations upon compression, we observe an isotropic (*s*-wave)  $\Delta$  with a large gap-to-superconducting transition temperature ratio,  $2\Delta_0/k_B T_c = 5.3(2)$  [ $\Delta_0 = \Delta(0\text{ K})$ ].  $2\Delta_0/k_B T_c$  decreases continuously upon pressurization until it approaches a value of  $\sim 3.5$ , characteristic of weak-coupling BCS theory of superconductivity despite the dome-shaped dependence of  $T_c$  on interfullerene separation. The results indicate the importance of the electronic correlations for the pairing interaction as the metal/superconductor-insulator boundary is approached.

The family of unconventional superconductors, that do not fit into the framework of the Bardeen-Cooper-Schrieffer (BCS) theory, has grown considerably over the last couple of decades and now includes cuprates<sup>1</sup>, heavy-fermions<sup>2</sup>, organic superconductors<sup>3</sup> and most recently also iron pnictides<sup>4</sup>. They all share a similar phase diagram<sup>5,6</sup> – superconductivity emerges through doping or applied pressure when the competing magnetic state is suppressed. Remarkably, even after more than 20 years of intensive research the superconducting pairing mechanism is still not fully understood for these compounds<sup>7</sup>. However, unlike in the phonon-driven BCS superconductors, strong electron correlations and magnetic interactions are believed to be important for the Cooper pairing mechanism.

A comparable phase diagram has been recently established for the cubic alkali fullerenes  $\text{A}_3\text{C}_{60}$  ( $\text{A}$  = alkali metal)<sup>8–10</sup> with the unit cell volume per fullerene ion,  $V$ , as a controlling parameter. For small  $V$ , short distances between neighbouring  $\text{C}_{60}^{3-}$  anions result in a strong overlap of the highest occupied triply-degenerate  $t_{1u}$  molecular orbitals and stabilize a Fermi-liquid metallic state from which the superconductivity emerges<sup>11,12</sup>.  $T_c$  at first increases with  $V$  but then for the optimal  $V^{\text{max}}$  it reaches a maximum  $T_c^{\text{max}} = 35$  and 38 K for the face-centred (f.c.c.) and A15 cubic polymorphs, respectively<sup>8–10</sup>. For even larger  $V$ , which are accessible only with the  $\text{Cs}_3\text{C}_{60}$  composition under pressure,  $T_c$  starts to decrease with increasing  $V$ . At the critical volume  $V_m$ , the on-site electron-electron Coulomb repulsion energy ( $U$ ) prevails over the electronic kinetic energy (measured by the bandwidth  $W$ ) and the metal/superconductor-to-Mott Jahn Teller-insulator transition (MIT) takes place. The importance of the Jahn-Teller effect arises from the intrinsic orbital degeneracy of the fullerenes<sup>13</sup>.

The size and symmetry of the superconducting gap,  $\Delta$ , characterize the superconducting state. In the case of cuprates, the *d*-wave symmetry of the superconducting gap<sup>14</sup> and its large amplitude ( $2\Delta_0/k_B T_c$  is much larger than the weak-coupling BCS value of 3.52<sup>15,16</sup>; here  $\Delta_0$  is the value of  $\Delta$  extrapolated to 0 K) are considered as fingerprints of an unconventional superconducting state. On the other hand, isotropic (*s*-wave) gaps with



**Figure 1 | Phase coexistence.**  $^{13}\text{C}$  (a) and  $^{133}\text{Cs}$  (b) NMR spectra of f.c.c.  $\text{Cs}_3\text{C}_{60}$  measured at 11 K in the superconducting state as a function of pressure. The  $^{13}\text{C}$  ( $^{133}\text{Cs}$ ) NMR spectra measured at ambient pressure in the paramagnetic Mott-insulating state are also shown for comparison. At low pressure ( $0.5 \leq P \leq 1.3$  kbar), two contributions to the spectrum from the superconducting (red lines) and Mott-insulating (blue lines) phases are present. The grey vertical line in (a) marks the expected (chemical) shift in the superconducting state.

substantially different sizes for different Fermi pockets are found in iron-pnictides<sup>17</sup>, where Hund's rules play an important role in controlling the electronic structure<sup>18</sup>. Despite these differences, the scaling of  $\Delta_0$  with  $T_c$  in the underdoped regime in both families<sup>16,17</sup> suggests that  $\Delta_0$  is controlled by the pairing strength. In the case of the extensively studied  $\text{A}_3\text{C}_{60}$  molecular superconductors with  $V < V^{\text{max}}$ , such as  $\text{K}_3\text{C}_{60}$  and  $\text{Rb}_3\text{C}_{60}$ , the isotropic  $\Delta$  found in numerous NMR or  $\mu\text{SR}$  experiments<sup>11,19–22</sup> has been for many years regarded as a result of a standard type-II BCS superconductivity in the weak coupling limit. However, the strong on-site electron-electron repulsion ( $U \sim 1$  eV)<sup>23</sup>, which is comparable with or even larger than the bandwidth ( $W \sim 0.5$  eV)<sup>11,12</sup>, suggests the importance of electron correlations and casts doubts on the applicability of the BCS formalism<sup>24</sup> thus bringing forward models of superconductivity which explicitly include correlations<sup>25–28</sup>. For  $V > V^{\text{max}}$ , the decrease of  $T_c$  with  $V$  found for both polymorphs of  $\text{Cs}_3\text{C}_{60}$  under pressure<sup>9,10</sup> is a strong indication of the growing importance of electron correlations. Although experiments on expanded  $\text{Cs}_3\text{C}_{60}$  have precisely determined the locations of the superconducting, normal and Mott-insulating states on the phase diagram<sup>9,10,29–31</sup>, the key information about the size and the symmetry of the superconducting gap in the vicinity of the parent Mott-Jahn-Teller insulating state is still missing.

In this work, we have used the local-probes  $^{13}\text{C}$  and  $^{133}\text{Cs}$  NMR at high hydrostatic pressures to study the f.c.c.  $\text{Cs}_3\text{C}_{60}$  polymorph as it is driven back from the ambient-pressure Mott-insulating state to a metallic/high- $T_c$  superconducting state by compression. We find that as the pressure is released and  $V$  increases, the decrease in  $T_c$  in the vicinity of the metal-insulator (MI) boundary is accompanied by the significant enhancement of the superconducting ratio,  $2\Delta_0/k_B T_c$  while the  $s$ -wave symmetry of the superconducting order parameter is retained for all  $V$ . The BCS theory fails to account for these results, which thus provide very stringent tests for the fundamental mechanisms that are responsible for superconductivity in f.c.c.  $\text{A}_3\text{C}_{60}$ .

## Results

The high f.c.c.  $\text{Cs}_3\text{C}_{60}$  (86%) phase fraction of the present sample has allowed us to use  $^{13}\text{C}$  and  $^{133}\text{Cs}$  NMR spectroscopy as a local probe of

the normal and superconducting state properties of the f.c.c.  $\text{Cs}_3\text{C}_{60}$  polymorph with applied pressure (Fig. 1, Fig. S1). At ambient pressure, the large paramagnetic susceptibility of the exchange-coupled Mott-insulating state at low temperature dramatically broadens the NMR spectra through the hyperfine interaction<sup>10</sup> and, for instance, prevents the clear separation of octahedral and tetrahedral  $^{133}\text{Cs}$  signals. The  $^{13}\text{C}$  NMR shift calculated from the first moment of the spectra, is shifted by about 190 ppm relative to the TMS standard, a characteristic value of  $\text{C}_{60}^{3-}$  ions (Fig. 1a). At the same time, the  $^{13}\text{C}$  spin-lattice relaxation rate (Fig. S2),  $1/T_1$ , is temperature independent consistent with the insulating nature of the electronic ground state down to 4 K<sup>10</sup>.

Slight pressurization of the sample leads to the appearance in the spectra of an additional much narrower component in addition to the broad NMR signal due to the paramagnetic insulator. In the case of the  $^{133}\text{Cs}$  NMR spectra measured at 0.5 and 1.3 kbar (Fig. 1b), this is seen as the emergence of a two-peak component with a peak intensity ratio of 2:1 – reflecting the relative occupancy by the  $\text{Cs}^+$  ions of the tetrahedral and octahedral sites<sup>32</sup>, respectively, in the crystal structure – superimposed on the broad signal. Similarly, the  $^{13}\text{C}$  NMR spectra show low-temperature narrowing (Fig. 1a) below approximately 25 K at 0.9 kbar. This narrowing of the NMR spectra with applied pressure provides the signature of the transition to the superconducting state in which the spin susceptibility vanishes, affording the sharper signals which come from the superconductor. The appearance of the superconducting component is also picked up very sensitively in the spin-lattice relaxation rate data (Fig. S2). For instance, slight pressurization ( $P = 0.9$  kbar) of the f.c.c.  $\text{Cs}_3\text{C}_{60}$  sample leads to an incomplete suppression of  $1/T_1$  below  $T_c = 25$  K that is consistent with the opening of the superconducting gap,  $\Delta$ . However, since  $1/T_1$  does not approach zero even when the temperature is reduced well below  $T_c$ , the transition here to the superconducting state is incomplete. The partial suppression of the spin-lattice relaxation rates and the presence of two overlapping NMR components are thus consistent with the presence of superconducting and Mott-insulating phases in different parts of the sample at these pressures. Such phase coexistence close to the MIT



boundary has also been detected for the A15  $\text{Cs}_3\text{C}_{60}$  polymorph<sup>9</sup> and implies that the transition between the Mott-insulating and superconducting states is of first order.

As the pressure is gradually increased, the narrow NMR signal due to the superconducting component completely dominates the low-temperature NMR spectra in agreement with the presence of bulk superconductivity for  $P \geq 1.7$  kbar (Fig. 1). The critical temperature has increased to 26.5(5) K at 1.7 kbar. At this pressure, the temperature dependence of the  $^{13}\text{C}$  NMR line shift, calculated from the first moment of the spectra, is rapidly suppressed for  $T \leq T_c$  (Fig. 2). We note that in the superconducting state there are three main contributions to the  $^{13}\text{C}$  NMR shift: the temperature-independent chemical shift, the Knight shift, which is proportional to the spin susceptibility, and the diamagnetic contribution due to the Meissner effect. The latter is estimated to be very small – around 3.2 ppm<sup>33</sup> – and can thus be neglected. The  $^{13}\text{C}$  NMR shift approaches  $\sim 150$  ppm as  $T \rightarrow 0$ , which is the value expected for the  $\text{C}_{60}^{3-}$  isotropic chemical shift<sup>33</sup>. We thus conclude that the Knight shift vanishes at  $T = 0$  thus providing firm evidence for the vanishing spin susceptibility of the spin-singlet Cooper pairs. The complete suppression of the  $^{13}\text{C}$  and  $^{133}\text{Cs}$  spin-lattice relaxation rates well below  $T_c$  also fully supports such a conclusion. However, there is no characteristic enhancement of the spin-lattice relaxation rates just below  $T_c$  that would mark the presence of the Hebel-Slichter coherence peak [Fig. 2b & d].

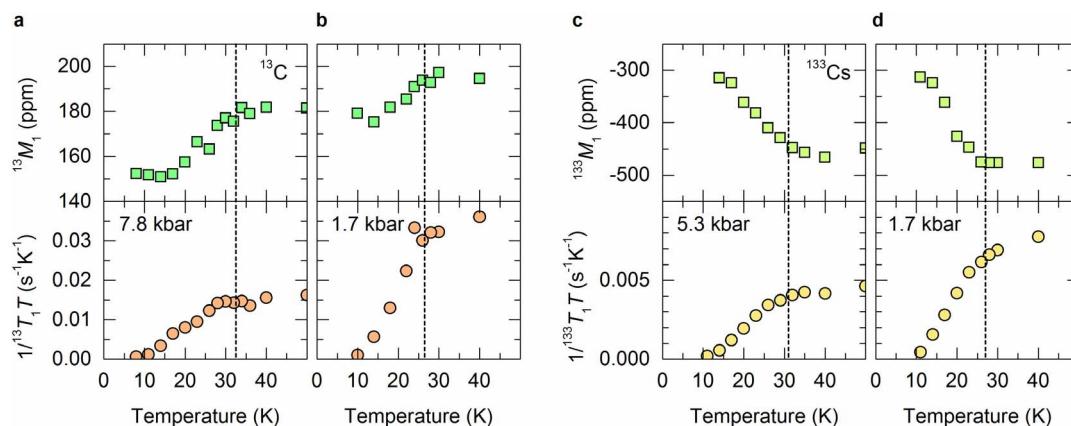
The temperature dependence of the spin-lattice relaxation rates below  $T_c$  can provide information on the size and symmetry of the superconducting gap,  $\Delta$ . Plotting  $(1/T_1)$  against  $1/T$  as a semilog plot for the 1.7 kbar data yields a straight line [Fig. 3a & Fig. S3]. This implies that the temperature dependence of the spin-lattice relaxation rate follows an activated behaviour with a single isotropic BCS-like (*s*-wave) superconducting gap,  $\Delta_0$  as described by the equation:  $1/T_1 \propto \exp[-\Delta_0/k_B T]$ , where  $k_B$  is the Boltzmann constant. This result rules out other singlet-pairing symmetries, such as that of a *d*-wave for which a power-law dependence  $1/T_1 \propto T^3$  is expected<sup>34</sup>. However, at the lowest temperatures, the  $T_1$  values are unambiguously longer by at least a factor of four than those expected in the BCS weak-coupling limit implying strong enhancement of the superconducting gap. Taking into account data for  $(T_c/T) > 1.25$ , the magnitude of  $\Delta_0$  at 1.7 kbar is found to be 6.0(2) meV. At the same time, the normalised gap value,  $2\Delta_0/k_B T_c = 5.3(2)$  is significantly enhanced relative to that expected for a weakly coupled BCS superconductor ( $2\Delta_0/k_B T_c = 3.52$ ). The single *s*-wave symmetry superconducting gap in f.c.c.  $\text{Cs}_3\text{C}_{60}$  contrasts with the behaviour of other

high-temperature superconducting families. The cuprates, which have a parent Mott insulating state like the fullerenes and are accepted as strongly correlated, universally show *d*-wave superconductivity<sup>14</sup>.  $\text{MgB}_2$ , which is not correlated and has multiple bands at the Fermi level that can be compared with the three electronically active  $t_{1u}$  orbitals in  $\text{Cs}_3\text{C}_{60}$ , is *s*-wave but displays multiple gaps<sup>35,36</sup>. Similarly, the iron pnictides, which are weakly to moderately correlated systems with multiple bands, also show multiple *s*-wave gap behaviour<sup>18</sup>.

Having established the fundamental behaviour of f.c.c.  $\text{Cs}_3\text{C}_{60}$  close to the metal/superconductor-to-insulator transition, we next address how the superconducting gap relates to the superconducting  $T_c$  upon tuning the bandwidth by pressure. As the applied pressure increases and the interfullerene separation gradually decreases,  $T_c$  (Fig. 3b) first increases (29(1) K at 2.9 kbar), reaches a maximum at 33(1) K at 7.8 kbar and then begins to decrease to 30.5(8) K at the highest pressure of the present experiments (14.2 kbar) tracking the dome-shaped,  $T_c(P)$  response established before by low-field magnetisation measurements<sup>10</sup>. However, contrary to this behaviour of  $T_c$ , the gap,  $\Delta_0$  does not show a maximum value but rather decreases monotonically (Fig. S3) with increasing bandwidth (decreasing  $V$ ). When the normalised value of the gap,  $2\Delta_0/k_B T_c = 4.7(2)$  (at 2.9 kbar) is considered (Fig. 3c), we find that this smoothly decreases towards the BCS weak coupling value of 3.52 at 7.8 kbar and above, implying a continuous reduction in the coupling strength as the system moves away from the metal-to-insulator boundary.

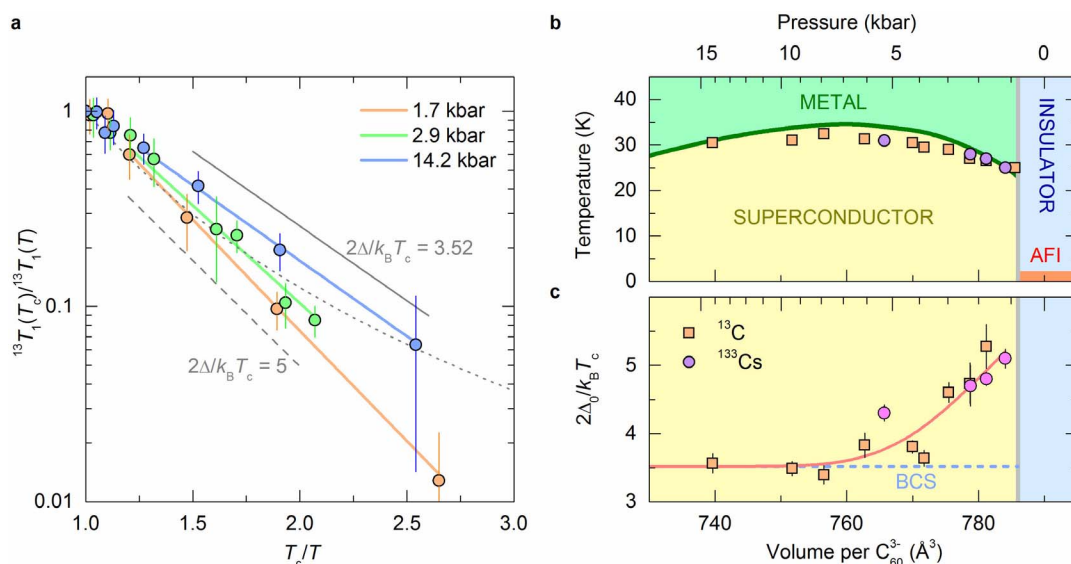
## Discussion

The enhancement of  $2\Delta_0/k_B T_c$  (Fig. 3c) with a concomitant decrease in  $T_c$  (Fig. 3b) for the large  $V$  region of the phase diagram provides an unprecedented opportunity to test the applicability of the BCS theory in alkali fullerenes. In principle, the maximal value,  $2\Delta_0/k_B T_c \approx 5$  close to the MIT boundary, could be obtained for strong electron-phonon coupling, but this would require optical or intermolecular phonons ( $\omega_{ph} \sim 100 \text{ cm}^{-1}$ )<sup>37</sup> to take part in the superconducting pairing mechanism. The weak-coupling BCS value of 3.52 found at high  $P$  (small  $V$ ) can only be obtained by the involvement of the intramolecular phonons ( $\omega_{ph} \sim 1000\text{--}1500 \text{ cm}^{-1}$ ) in the pairing interaction. The  $V$  dependence of  $2\Delta_0/k_B T_c$  would thus then require that phonon modes in distinctly different spectral regions are active in different parts of the electronic phase diagram. This is unlikely as the intramolecular phonon modes are always present and cannot be active only in one part of the phase diagram. Therefore, these



**Figure 2 | Critical temperatures.** (a), (b) Temperature dependence of the  $^{13}\text{C}$  NMR shifts (green squares) and the spin-lattice relaxation rates,  $1/T_1 T$  (orange circles) measured at 7.8 and 1.7 kbar, respectively. (c), (d) Temperature dependence of the  $^{133}\text{Cs}$  NMR shifts (light green squares) and the spin-lattice relaxation rates,  $1/T_1 T$  (light orange circles) measured at 5.3 and 1.7 kbar, respectively. The  $^{13}\text{C}$  ( $^{133}\text{Cs}$ ) NMR shifts were obtained from the first moments,  $M_1$ , of the  $^{13}\text{C}$  ( $^{133}\text{Cs}$ ) NMR spectra. The dashed vertical lines mark the onset temperatures at which  $M_1$  becomes suppressed in the superconducting state. At high pressure [(a), (c)],  $1/T_1 T$  is suppressed at a slightly lower temperature than  $M_1$ , which is the signature of a damped Hebel-Slichter coherence peak. At low pressure [(b), (d)], the two onset temperatures coincide, thus implying the absence of a coherence peak.





**Figure 3 | Superconducting gap.** (a)  $^{13}\text{C}$  spin-lattice relaxation rate,  $1/T_1$  normalized to its value at  $T_c$  vs inverse temperature,  $T_c/T$  for three characteristic pressures 1.7 (orange circles), 2.9 (green circles), and 14.2 kbar (blue circles). Solid lines through the points are fits to the equation,  $1/T_1(T) = A \exp[-\Delta_0/k_B T]$ , where the fitting parameters are the amplitude  $A$  and the value of the superconducting gap at  $T = 0$  K,  $\Delta_0$ . Only data for  $(T_c/T) > 1.25$  are included in the fits. Thin solid and dashed lines mark the expected slopes for  $2\Delta_0/k_B T_c$  ratios of 3.52 and 5, respectively. The dot-dashed line is the power law dependence,  $1/T_1 \propto T^3$  anticipated for  $d$ -wave superconductivity. (b) The low-temperature phase diagram of f.c.c.  $\text{Cs}_3\text{C}_{60}$  as derived from the shifts of the NMR spectra in the superconducting state. Squares and circles mark the onset of superconductivity as deduced from the  $^{13}\text{C}$  and  $^{133}\text{Cs}$  NMR data, respectively. The experiments were conducted at a 9.39 T magnetic field. The volume (pressure, top scale) dependence of  $T_c$  represented by the solid green line is that obtained from the low-field magnetization studies on the same sample<sup>10</sup>. The thick grey vertical line indicates the critical volume,  $V_m$  for the metal/superconductor-to-Mott insulator transition. The antiferromagnetic transition temperature,  $T_N = 2.2$  K, of the Mott-insulating phase is taken from Ref. [10] (AFI denotes antiferromagnetic insulating phase). (c) The volume per  $\text{C}_{60}^{3-}$ ,  $V$ , dependence of the superconducting gap divided by the superconducting critical temperature,  $2\Delta_0/k_B T_c$ , obtained from the  $^{13}\text{C}$  (squares) and  $^{133}\text{Cs}$  (circles) spin-lattice relaxation rate data in the superconducting state. The solid thick line is a guide to the eye, while the dashed blue line marks the BCS value,  $2\Delta_0/k_B T_c = 3.52$ . The thick grey vertical line marks the metal/superconductor-to-Mott insulator critical volume,  $V_m$ .

arguments rule out the conventional BCS-like explanation of superconductivity in f.c.c.  $\text{Cs}_3\text{C}_{60}$ , despite the retention of  $s$ -wave symmetry over the entire phase diagram.

The failure of the BCS theory therefore necessitates the presence of an additional parameter responsible for the strong coupling superconductivity as we approach the MIT boundary from the low volume per  $\text{C}_{60}^{3-}$ ,  $V$ , side of the phase diagram. In this part of the phase diagram, the screening of the Coulomb interactions is not so effective anymore and the effective  $U$  is expected to rapidly increase thus making electron correlations progressively more important. For instance, they are directly reflected in the non-Korringa temperature dependence of  $1/T_1 T$  in the normal state for large  $V$ <sup>31</sup>. Therefore, we propose that the additional parameter in the superconducting mechanism is most likely the increased importance of correlations relative to the electronic bandwidth.

The  $s$ -wave nature of the superconducting gap and the anomalously large values of  $2\Delta_0/k_B T_c$  set the fullerenes out as an unusual class of strongly correlated superconductors in which the  $s$ -wave nature is retained right across the entire  $V$  range but the non-BCS nature is strongly indicated by the  $V$  (or bandwidth) dependence of the superconducting gap. This conclusion is in qualitative agreement with the recent application of density functional theory for superconductors within the local density approximation (LDA) to the alkali fullerenes,  $\text{A}_3\text{C}_{60}$  ( $\text{A} = \text{K}, \text{Rb}, \text{Cs}$ )<sup>24</sup>. This work strongly suggests the necessity to go beyond the framework of the Migdal-Eliashberg theory and calls for direct comparisons with the other families of high-temperature superconductors to determine whether they share the same purely electronic pairing mechanism. In the strongly correlated cuprates, the superconducting gap has a  $d$ -wave symmetry in momentum space with the gap values at the “antinodes” that place  $2\Delta_0/k_B T_c$  well in the strong coupling regime<sup>15</sup>. The gap generally first

increases with doping in the underdoped regime, varying in a similar manner to  $T_c$ , but then it soon reaches a maximum and remains at the same size for larger doping levels<sup>16</sup>. On the other hand, the pnictides are less correlated systems, but like the fullerenes they are multiband superconductors. Their superconducting gaps also scale linearly with  $T_c$  in the underdoped regime<sup>17</sup> and show many similarities with the behaviour observed in underdoped cuprates. The opposing volume dependences of  $2\Delta_0/k_B T_c$  and  $T_c$  observed for the f.c.c.  $\text{Cs}_3\text{C}_{60}$  compound are thus qualitatively different from these two families. Finally, we note that a preprint recently appeared on the arXiv server<sup>38</sup> reporting NMR results on the  $\text{A15 Cs}_3\text{C}_{60}$  polymorph that show a comparable  $V$  dependence of the superconducting gap thus implying a lattice-independent superconductivity mechanism in both these systems.

## Conclusions

The superconducting gap size in f.c.c.  $\text{Cs}_3\text{C}_{60}$  increases as the insulator is approached, but the singlet  $s$ -wave nature of the superconductivity is maintained across the entire phase diagram. The observation of a single isotropic gap from a molecular  $s/p$  system with three degenerate orbitals contributing to the states at the Fermi energy contrasts with the multiband multigap structure of extreme BCS  $\text{MgB}_2$ , which shares the electronic orbital parentage of the states involved but with much broader bands due to its extended non-molecular nature. The weaker overlap between the parent orbitals in the fullerenes and the observation of magnetism and electronic correlations in the competing states makes the contrasting variation of  $\Delta_0$  and  $T_c$  from the cuprates and the pnictides important. Hund’s rules are not relevant for the cuprates, as there is only one accessible spin state with one hole in the parent insulator, but are thought to control the electronic structure of the pnictides<sup>20</sup>. In the fullerenes,



there is a competition between high- and low-spin states involving both Hund's rules and Jahn-Teller electron-phonon coupling, which is decisive in favouring the low-spin state<sup>13</sup>. The observation of a large non-BCS but s-wave gap in a superconductor which emerges from an onsite correlation-driven Mott insulating state and in which the molecular nature of the electronic states gives strong local Jahn-Teller coupling may indicate that the fullerides are indeed non-BCS but are distinct from other examples of unconventional superconductivity in correlated materials.

## Methods

**Sample.** All measurements were performed on a Cs<sub>3</sub>C<sub>60</sub> sample that contains predominately f.c.c. (86%) polymorph with secondary A15 Cs<sub>3</sub>C<sub>60</sub> (3%), body-centred orthorhombic Cs<sub>4</sub>C<sub>60</sub> (7%) and CsC<sub>60</sub> (4%) phases, based on Rietveld refinement of synchrotron powder XRD data<sup>10</sup>.

**NMR measurements.** <sup>13</sup>C (nuclear spin  $I = 1/2$ ) and <sup>133</sup>Cs (nuclear spin  $I = 7/2$ ) NMR spectra and their spin-lattice relaxation times,  $T_1$ , were measured between 4 and 300 K at a magnetic field of 9.39 T. As references, tetramethylsilane (TMS) and CsNO<sub>3</sub> standards were used with corresponding reference frequencies,  $\nu(^{13}\text{C}) = 100.5713$  MHz and  $\nu(^{133}\text{Cs}) = 52.4609$  MHz, respectively. In the <sup>13</sup>C NMR lineshape measurements, a Hahn-echo pulse sequence  $\pi/2 - \tau - \pi - \tau - \text{echo}$  was used, with pulse length  $t_w(\pi/2) = 8$   $\mu\text{s}$  and interpulse delay  $\tau = 40$   $\mu\text{s}$ . In the <sup>133</sup>Cs NMR experiments, a two-pulse solid-echo sequence  $\pi/2 - \tau - \pi/2 - \tau - \text{echo}$  was used, with a pulse length  $t_w(\pi/2) = 5$   $\mu\text{s}$  and an interpulse delay  $\tau = 50$   $\mu\text{s}$ . The pulse length was optimized for <sup>133</sup>Cs nuclei in the high-symmetry octahedral and tetrahedral sites with zero quadrupole frequency (expected for the f.c.c. Cs<sub>3</sub>C<sub>60</sub> polymorph), which further suppressed the already weak signals of minority phases. For  $T_1$  measurements, the saturation recovery and inversion recovery techniques were both applied.

**High-pressure cell.** A home-built clamp-type cell design was used for the high-pressure NMR experiments. The cell body was manufactured from non-magnetic Ni-Cr-Al hardened alloy and 1:1 mixture of fluorinert oils FC-770 and FC-70 was used as pressure medium in order to minimize non-hydrostatic conditions. The ruby N<sub>2</sub> luminescence line was measured *in-situ* in order to follow the evolution of pressure inside the cell at each temperature. The unit-cell volume at each temperature and pressure was calculated from the published f.c.c. Cs<sub>3</sub>C<sub>60</sub> structural, thermal contraction and compressibility data<sup>10</sup>. Since the *in-situ* pressure was notably changing during the temperature dependent measurements, all reported pressures are given as those measured at 35 K. The absolute values of pressures are accurate within  $\pm 0.5$  kbar.

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## Author contributions

K.P., M.J.R. and D.A. designed and supervised the project. The samples were synthesized by A.G. and Y.T., who also performed the synchrotron X-ray diffraction and magnetic measurements, and analyzed the data. The NMR experiments were conducted and analyzed by A.P., A.K. and P.J. All authors contributed to the interpretation of the data and to the writing of the manuscript.

## Additional information

**Supplementary information** accompanies this paper at <http://www.nature.com/scientificreports>

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